

HBD clusterizer with built in background subtraction

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HBD meeting
2010.06.02

Intro 1

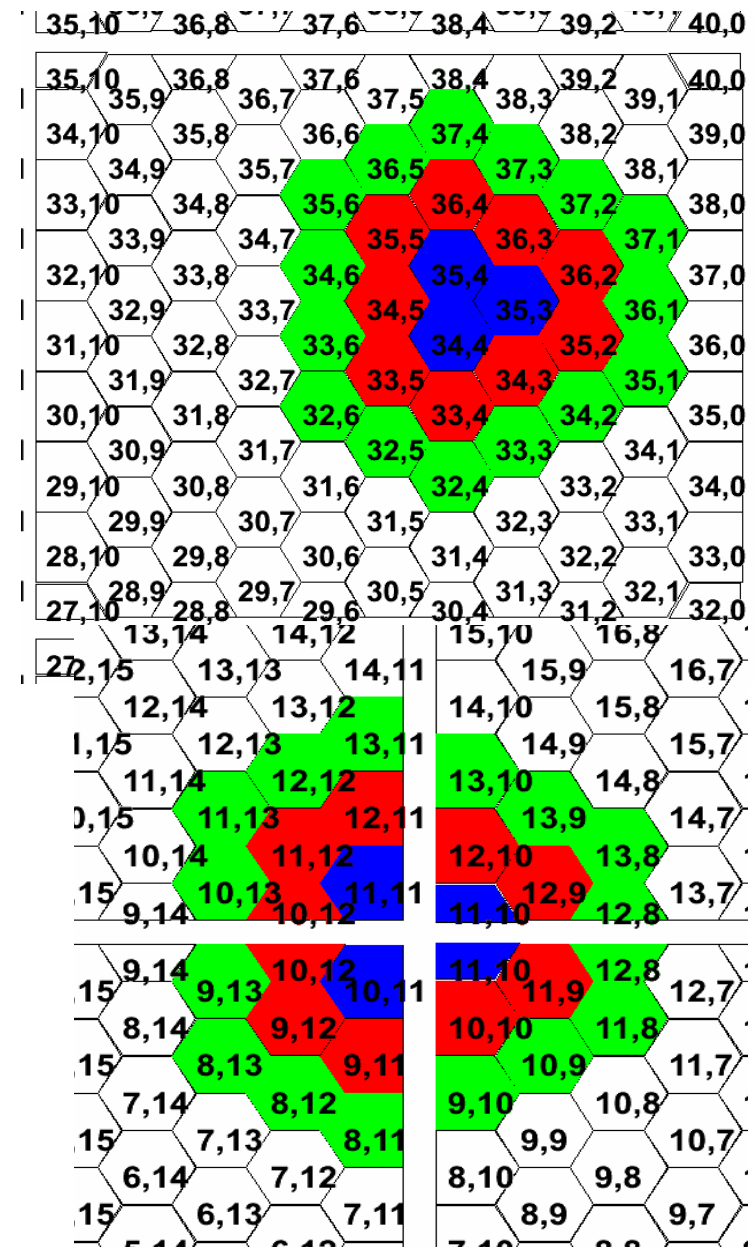
- An alternate clusterization algorithm is being developed
 - Main point: Background is handled by subtracting average per pad background estimated from surrounding area
 - It is still under development and testing but for people interested, its been submitted to cvs (in offline/analysis/hbd_proto)
 - It works like any other analysis module
 - Clusterizer: offline/analysis/hbd_proto/HbdLbsClusterizer
 - For this to work, one has to locally compile offline/packages/hbd after editing Makefile.am to add HbdBlobListv1.h to install headers
 - Embedding tests: offline/analysis/hbd_proto/HbdEmbed
 - Ntuples and plotting: offline/analysis/hbd_proto/HbdAnalysis
 - Simulation tuning: offline/analysis/hbd_proto/HbdMcChargeRecal
 - There is still a lot of debugging couts and some valgrind errors.
 - The code can be improved in efficiency and style
 - Any input is welcome, and feel free to modify if you have ideas or let me know

Intro 2: A new clusterization algorithm

- Better of the two worlds:
 - Like Weizmann clusterizer: two steps, “preclusterization” and merging.
 - But, before merging there is a control step where preclusters are selected based on a few criteria
 - Like HnS clusterizer: preclusters are triplets, most natural shape for the hexagonal symmetry of the HBD pads
 - It doesn't need to depend on the projection of electrons even in high background environment. Though this information can be used if needed.
- And a little bit more....
 - At the preclusterization step, a local background subtraction is internally (without the use of parametrization) applied.
 - This is done by estimating the background level from neighboring pads of the precluster. There seem to be (cf slide 5) reasonable correlation to warrant this
 - After merging, the final cluster's background is subtracted using neighboring pads
 - For this reason, will refer to the new clusterizer as of LBS (local background subtraction) method

Preclusterization

- First step of the algorithm is the selection of preclusters.
 - Candidates for preclusters are all possible compact triplets in the HBD (def. All members sharing a single edge with the other two members)
- Preclusters have
 - first neighbors
 - and second neighbors.
 - And they cross borders
- They have the following properties:
 - Charge & area of Members
 - Charge & area of 1st & 2nd neighbors
 - Net signal in the “member” zone
 - “Shape” meaning distribution of net charge among pads in member zone



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Justification of background estimation

- Basic assumption of the method
 - Scintillation background varies continuously over HBD surface
 - Background in any compact group of pads can be estimated from the average rate of npe in its neighboring pads

$$bkg = a_{mem} * \left(\frac{w_{fn} * q_{fn}}{a_{fn}} + \frac{(1 - w_{fn}) * q_{sn}}{a_{sn}} \right)$$

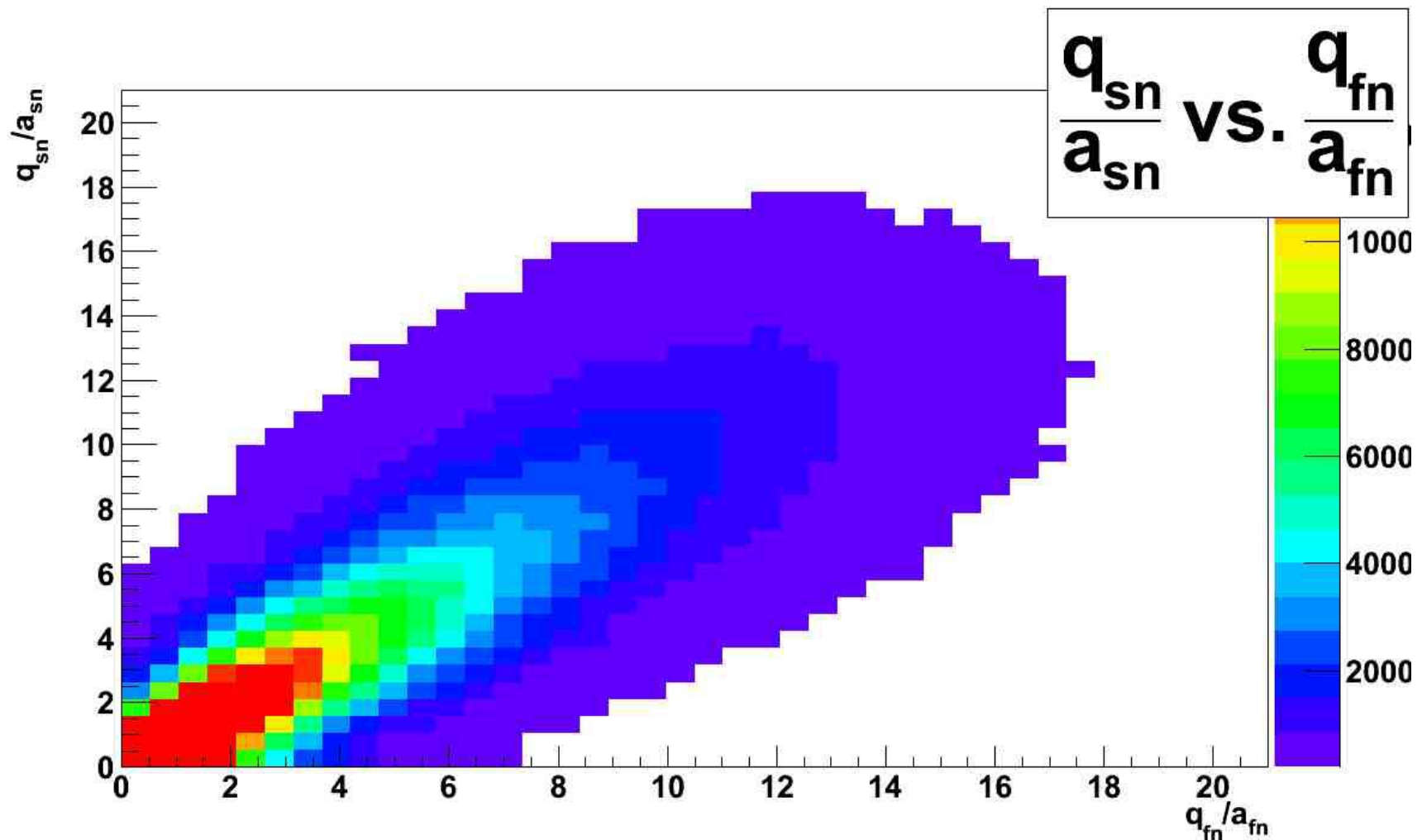
mem=triplet member fn=first neighbor, sn=second neighbor

a=area, q=number of photoelectrons

w= weight, for now set to 0.5

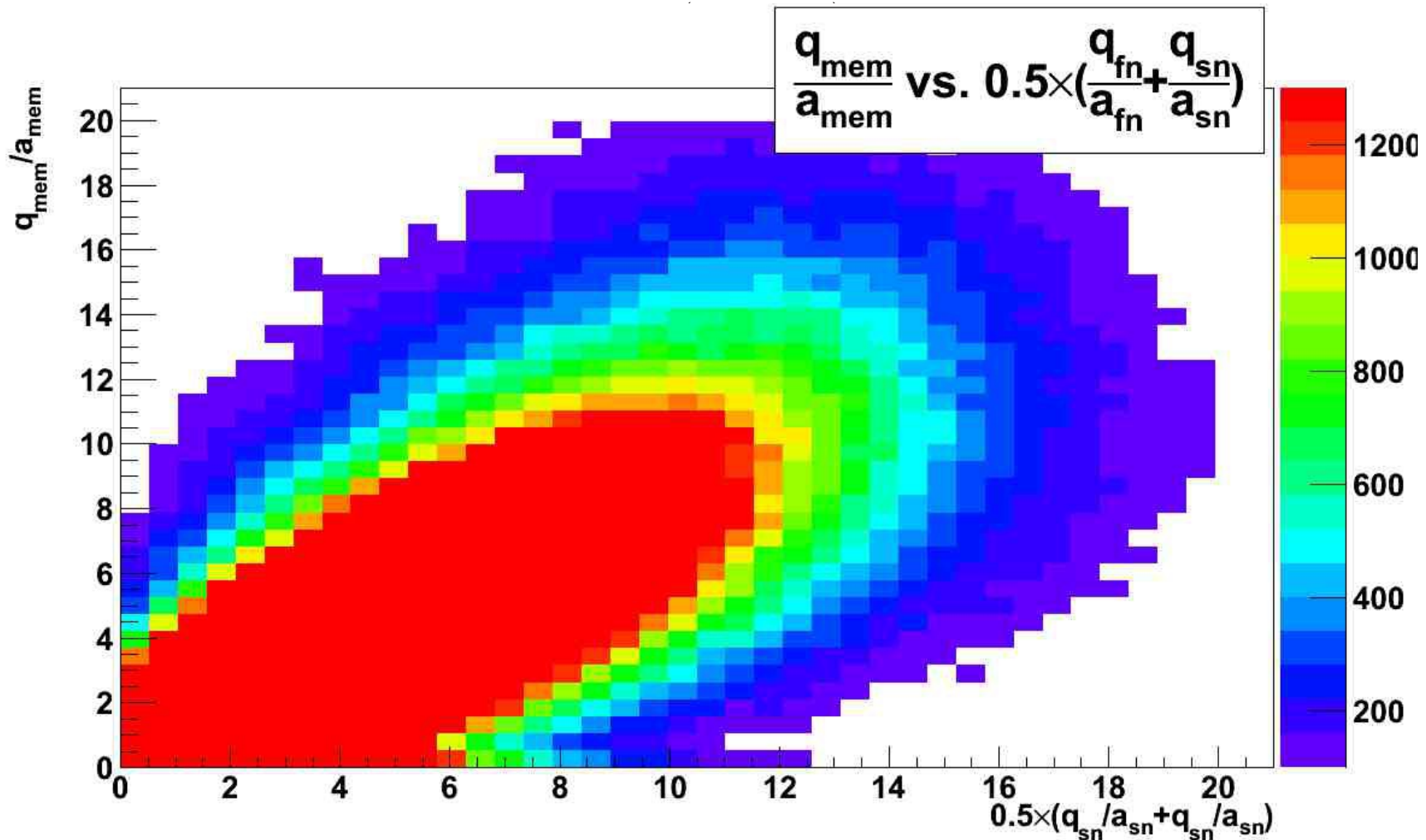
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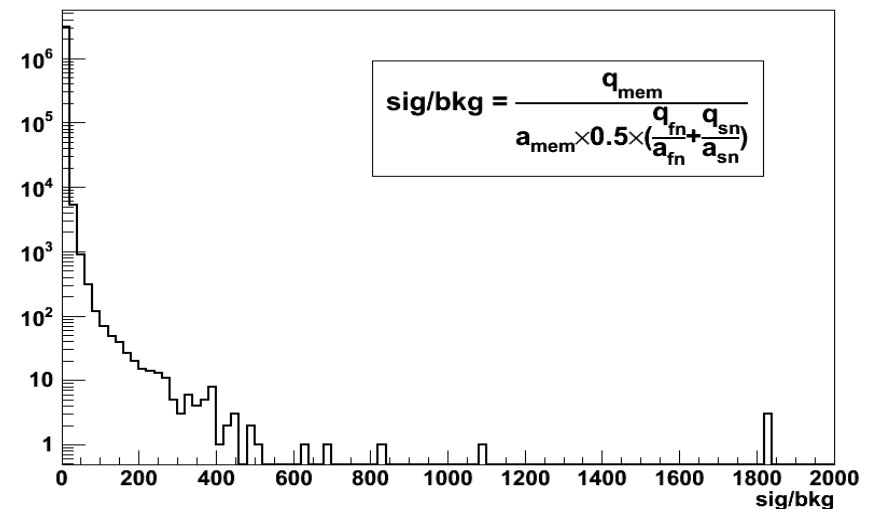
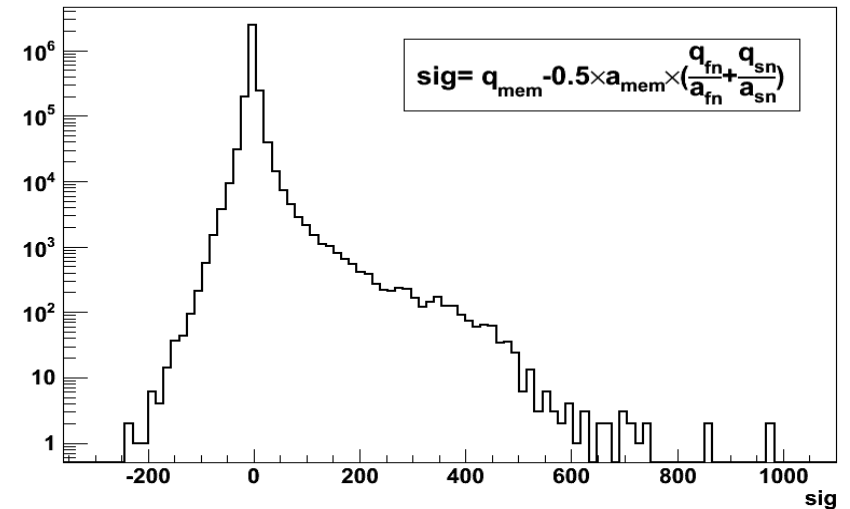
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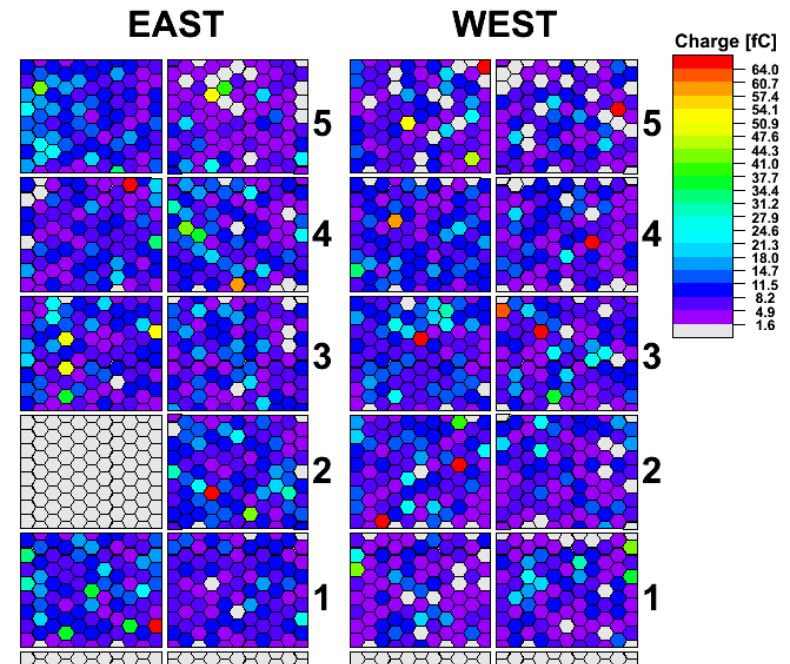
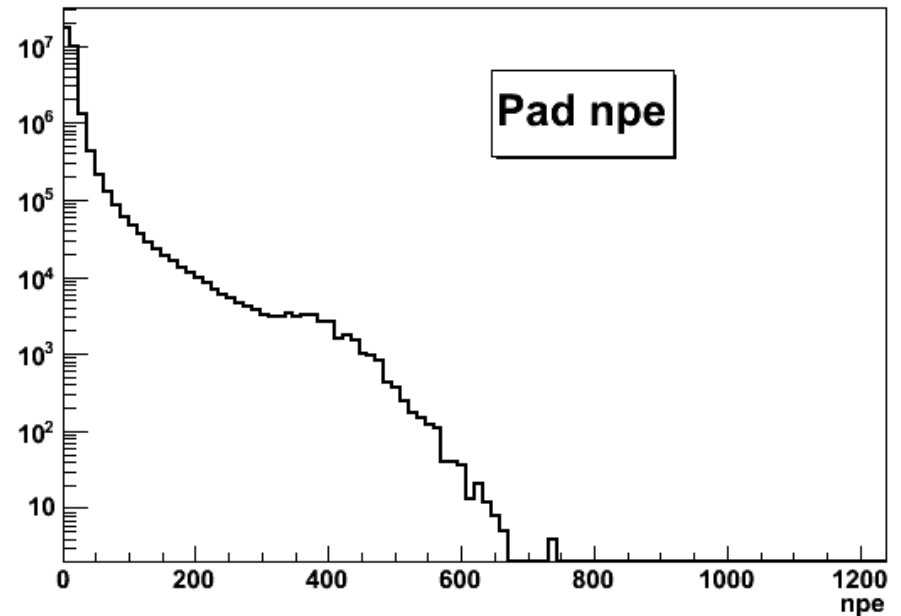
Precluster selection

- Don't want to keep everybody
 - Code will be slow
 - Will end up with superbig clusters
- What to keep?
 - Reasonable net signal
 - For now keeping $5 < \text{sig}(\text{npe}) < 50$
 - This spans both the singles and doubles expected charge in a triplet
 - Reasonable S/B
 - We can cut on estimated S/B
 - Optimization will be shown later
 - Shape cut
 - Distribution of a couple of such parameters will be shown later for data and MC



HIPs: an issue with a solution

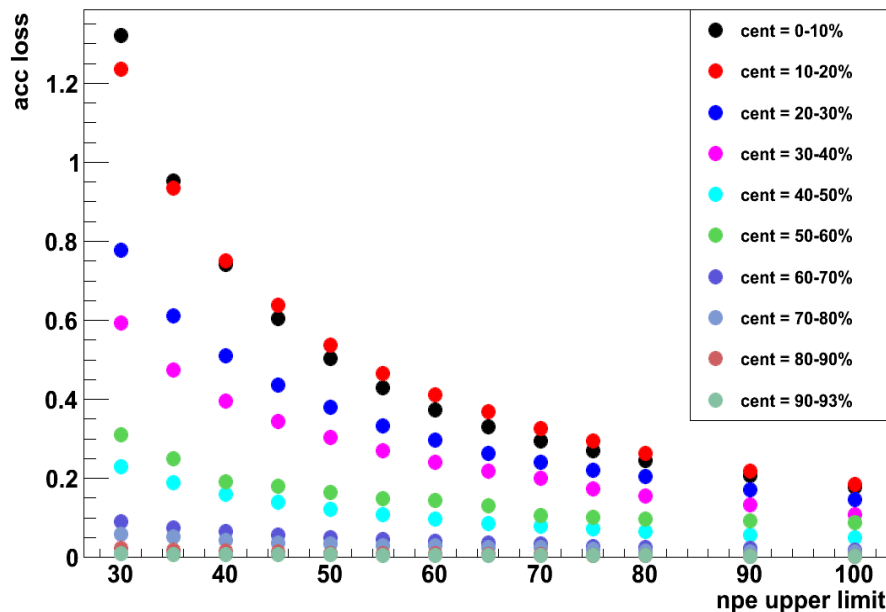
- The pad by pad charge distribution has a very long tail
 - Caused by physics processes that deposit a huge amount of energy
 - Much more than typical per pad charge expected from either scintillation or Cerenkov
 - Rate is proportional to intensity
 - X-ray, neutrons heavy particles?
- These pads if left alone are a big problem for any clusterization algorithm, because they can seed fake clusters.
- Fortunately, event by event, they cover only a very small fraction of the active HBD area



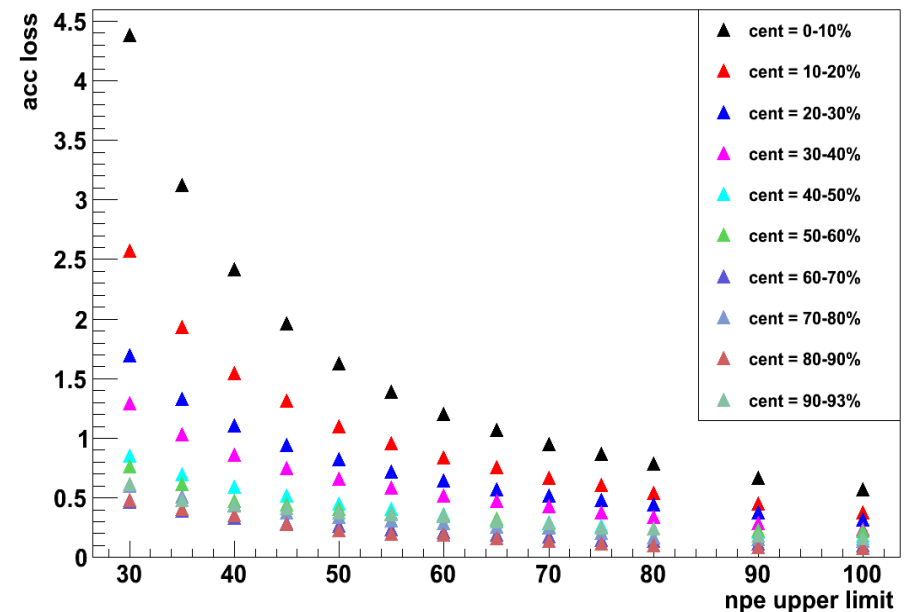
Effect of upper limit on pad npe

- Before clusterization one can set $npe=0$ for those pads that fire above a certain upper limit
 - Plot on left: Event averaged fraction of acceptance loss incurred by throwing out pads firing above an UL, vs. the value of the ul for different centralities
 - Plot on right: Fraction of pads firing above upper limit to those firing below upper limit but still above threshold

Fractional (%age) acceptance loss vs. per pad npe upper limit value



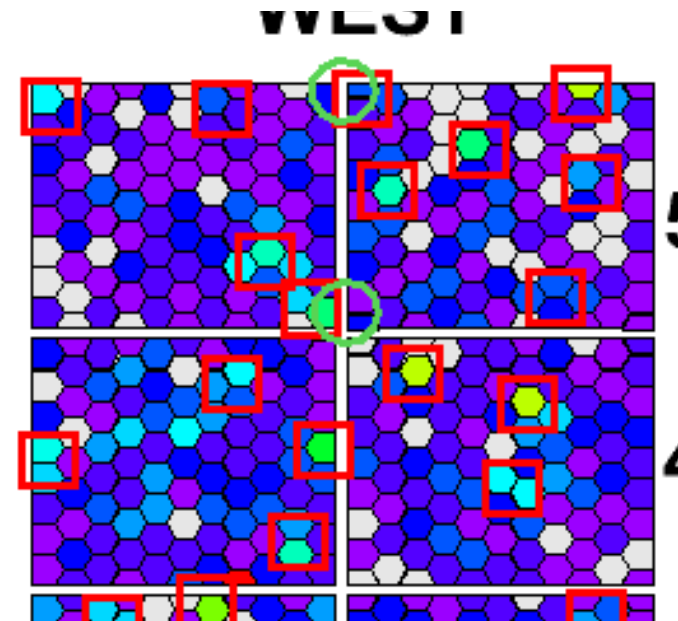
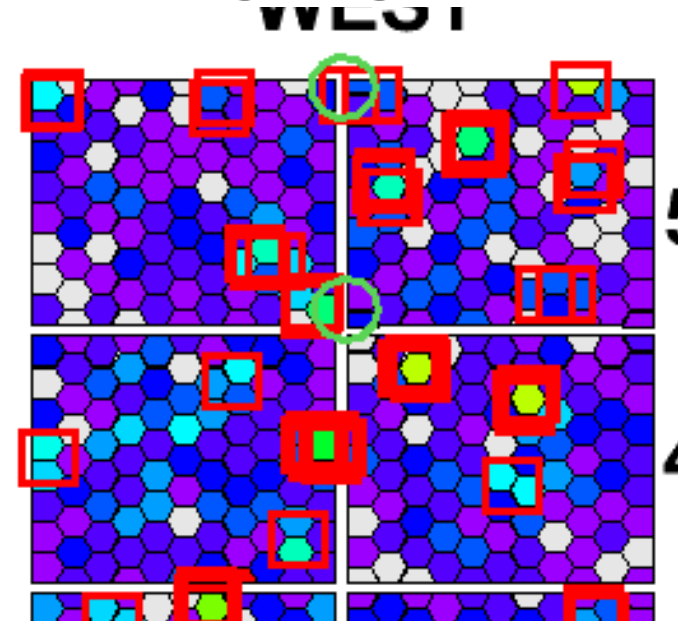
Fraction (%age) of hot pads to fired pads below upper limit vs. per pad npe upper limit value



Cutting at 50 seems safe. <2% of fired pads are lost even in most central event

Merging and post merging

- Overlapping preclusters
 - Share at least one pad
- Final clusters
 - Lump together pads from all overlapping groups of preclusters
- Local bkg. subtraction
 - Merged clusters have 1st and 2nd neighbors just like preclusters
 - 1st and 2nd neighbor charge is used to estimate background to subtract from the members of merged cluster
- Cluster track association
 - Nothing new here, based on proximity just like in Wis & HnS



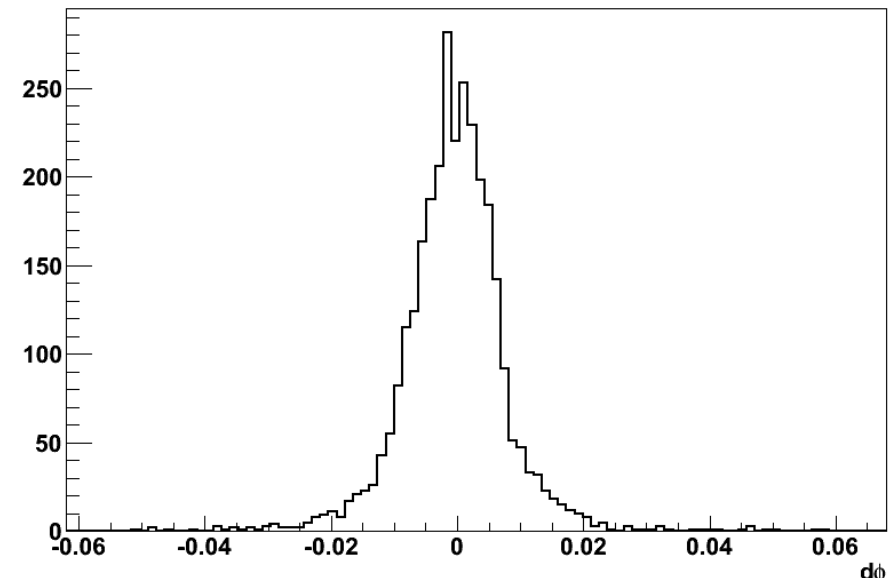
Validation

- For the validation here is the program
 - Single electron simulation with no background
 - Simulation tuning, geometry cross check, shape study
 - Single electron simulation with “emulated” background
 - Optimization of precluster selection criteria, fake rate, cluster size, cluster rates
 - Double electron simulation (Conv. and Dalitz) w/ & w/o emulated Bkg
 - Confirm doubling of the cluster signal, estimate misidentification rate from doubles created midway inside the HBD
 - p+p events
 - Event Accumulator/ Embedding
 - More realistic background. Do we still get same answer from the clusterizer for simulated electrons?
 - Embedding already implemented (initial test on single electrons)
 - Real Au+Au data
 - Cluster shape, singles/doubles/hadron charge comparison, Analysis

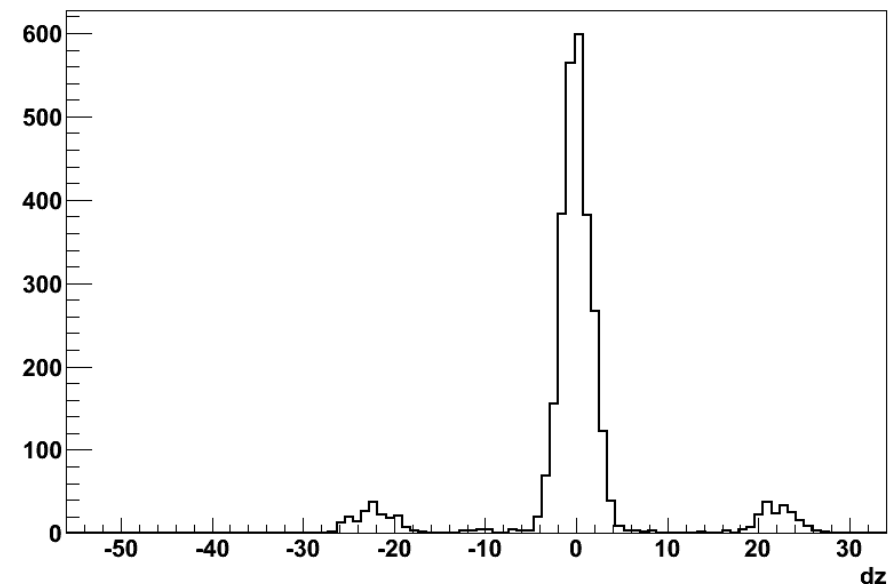
Single electrons, no background

- Usual PHENIX chain
 - $x, y, z = 0, 0, \pm 20$
 - Full Hbd response
- Run clusterizer
 - Dphi, Dz look very good
 - Except for wings at $\pm 20\text{cm}$ for dz
- This demonstrates that the geometry is being used correctly in the code.

dφ distribution for simulated electrons



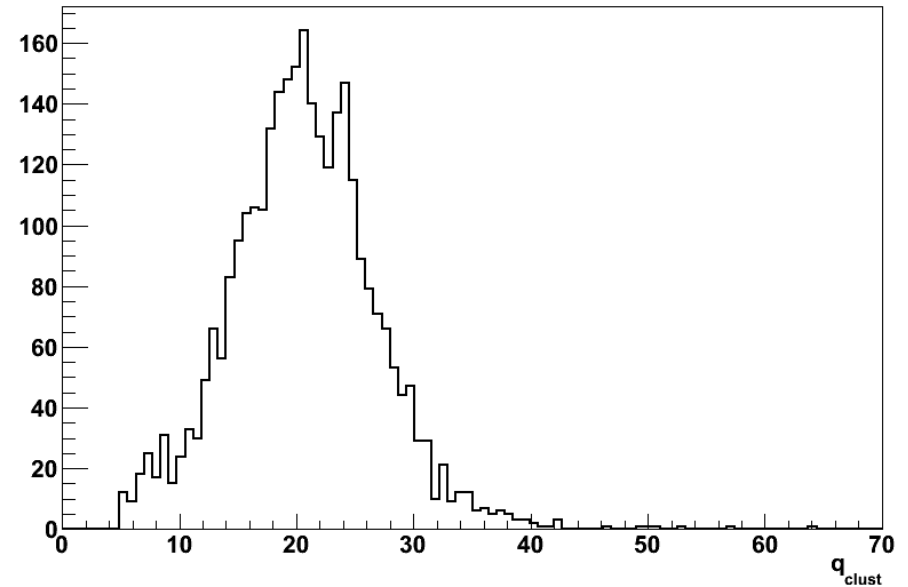
dz distribution for simulated electrons



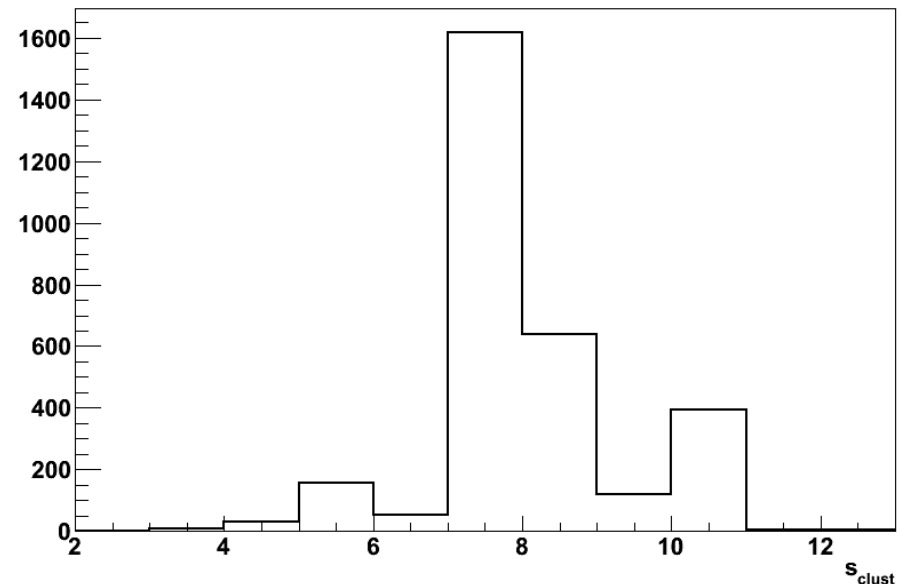
Cluster charge and size distributions

- Cluster charge distrib.
 - Off by a factor of ~ 2.7
 - Running a 'Recal' module that divides every pad by this factor
 - This should be done only on Cerenkov signal
- Cluster size distribution
 - Cluster sizes are somewhat big.
 - Current merging mechanism tends to add 1st neighbors
 - This should not have too much effect on the cluster charge since background is subtracted event by event

Cluster charge (npe) distribution for simulated electrons



Cluster size distribution for simulated electrons

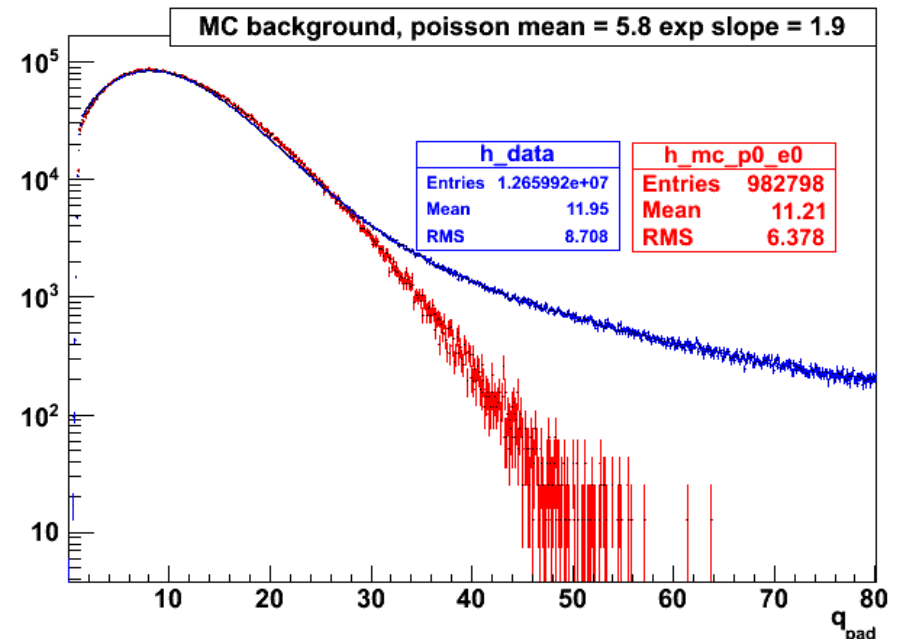
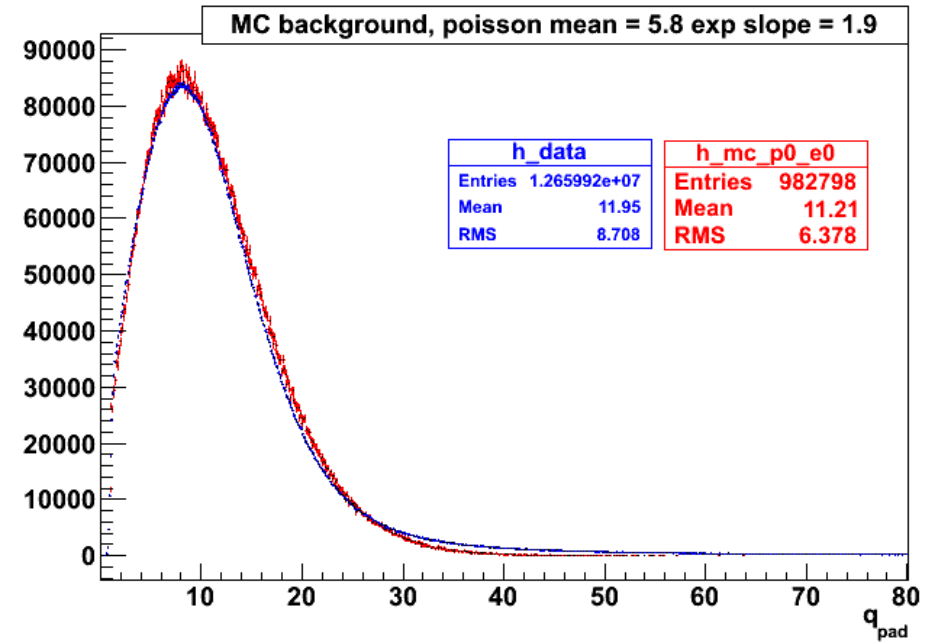


Mimic the real data background

- Attempt to generate RD like background

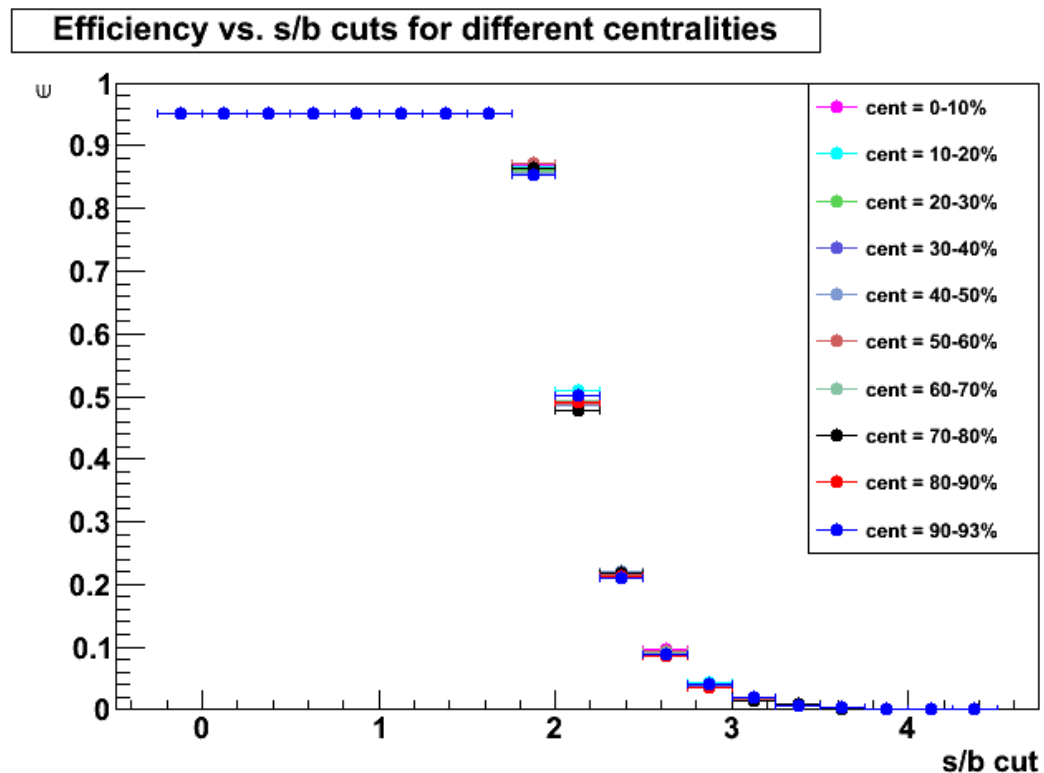
$$q = \sum_0^{P(M)} \exp(\tau)$$

- M (Poisson RV mean) and tau (Exp. RV decay const.) are hand tuned to match the RD pad charge distribution
 - Ten centrality bins of 10%
 - The long tail in RD is hard to reproduce (probably coming from jets? If so maybe can be added with some effort.)
 - This kind of detail matters for clusterizing
- Using temporarily as a rough approximation to scintillation background



Optimizing s/b precluster selection cut

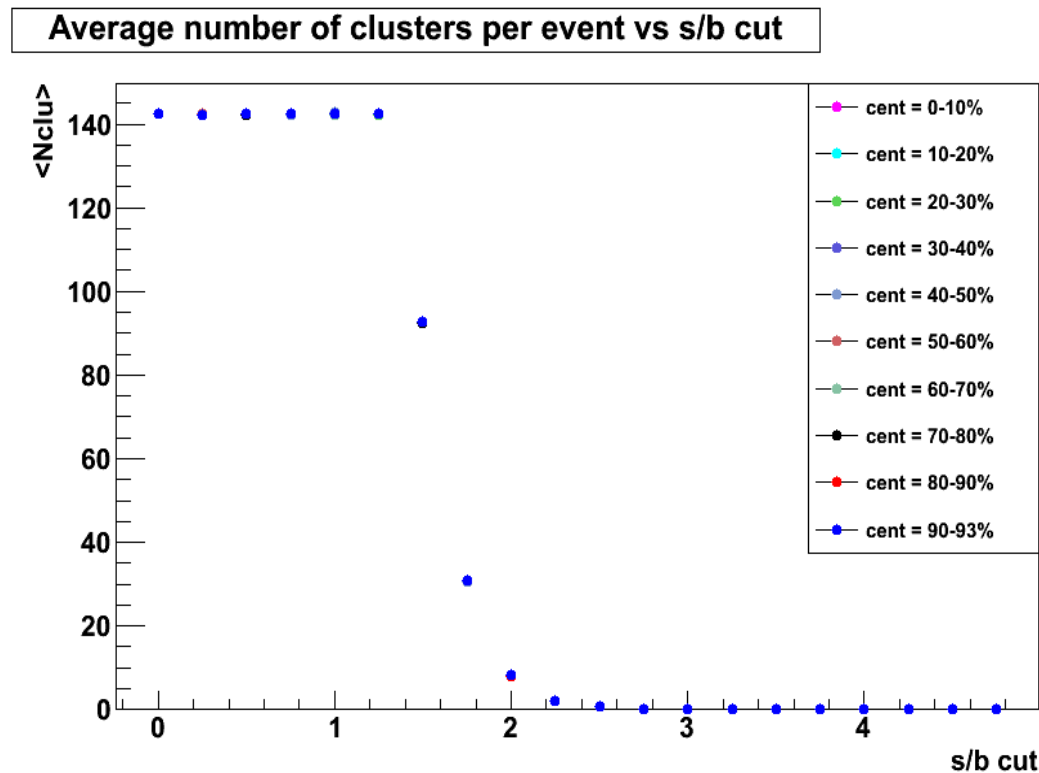
- Single electron cluster efficiency vs. s/b cut
 - Fraction of simulated single electrons that get associated with a cluster
 - vs. s/b cut using 'faux' scintillation background tuned to different centrality selections



- The sudden drop in efficiency happens at the same position for all centralities, which points to a possible problem with the background emulator
- Accumulator or embedding should give a better picture

Average number of clusters per event

- Similar structure as for the efficiency
 - The drop in number of clusters occurs earlier than for the efficiency
 - With a s/b cut at 2, the efficiency is still $> 90\%$ but $\langle N_{clu} \rangle$ is down to less than 10



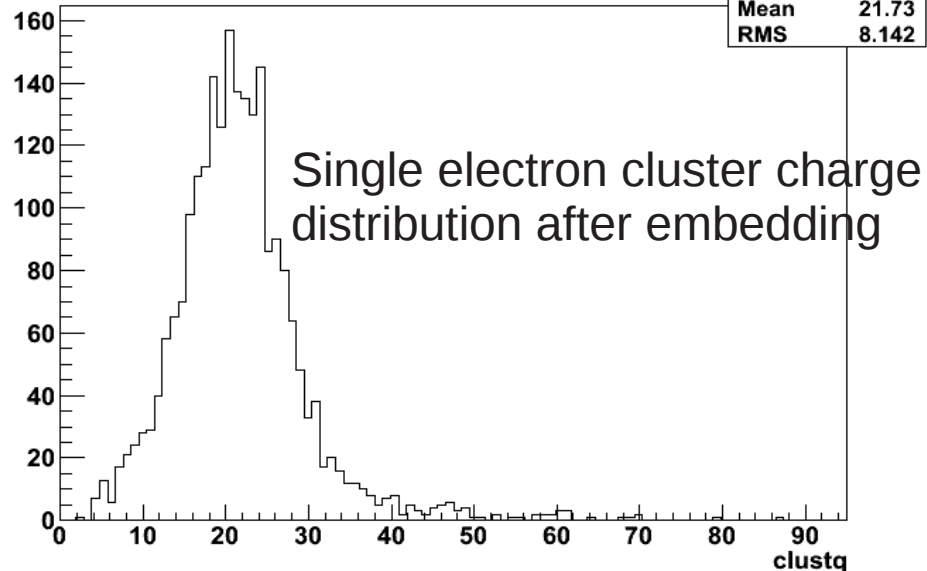
- This has to be confirmed by more realistic background env.
 - Embedding the simulated electron into events where there is no identified electron

Embedding MC Cerenkov response in RD

- Embedding is another option to see the effectiveness of a clusterization algorithm with real background
 - Simulate single (or double) electrons
 - Pad by pad add the signal from events in real data to the Cerenkov response from simulation
 - Run the clusterizer on merged HbdCellList
- Easy to implement real data event selection based on any criteria (bbcz, presence of electrons etc..) but not implemented yet. More to come...
- Embedding can be a useful tool to study the performance of a clusterizer (efficiency and stability in high background environment)

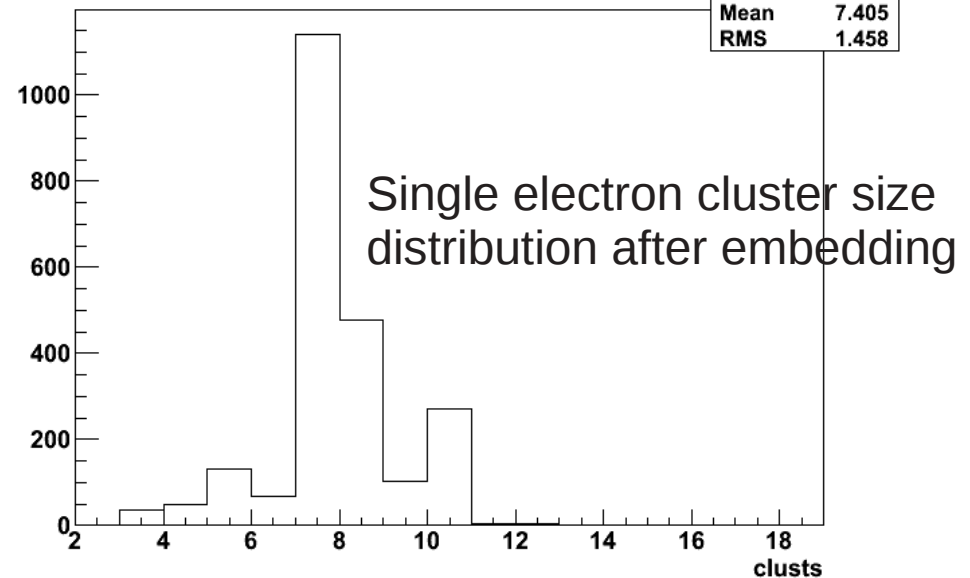
clustq {dphi!=-9999&&dz!=-9999}

htemp	
Entries	2285
Mean	21.73
RMS	8.142



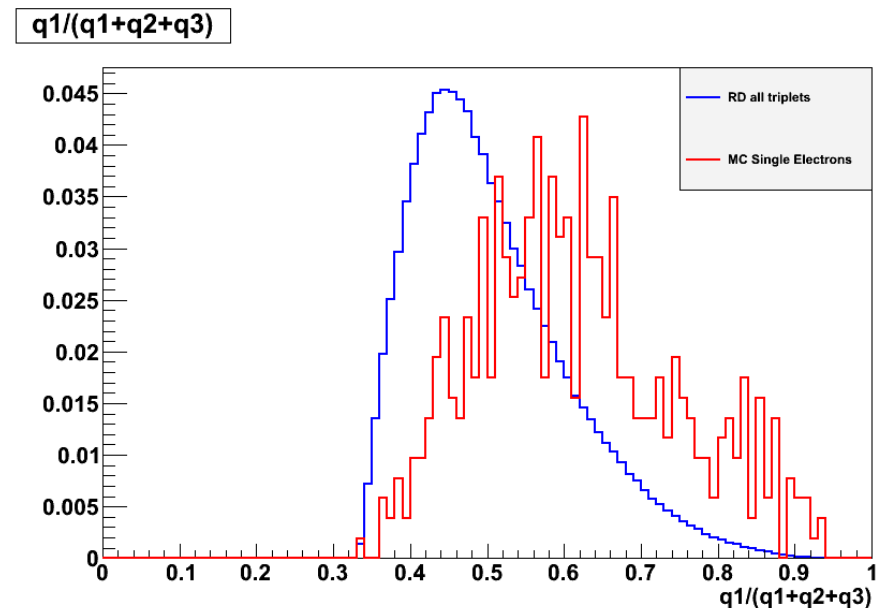
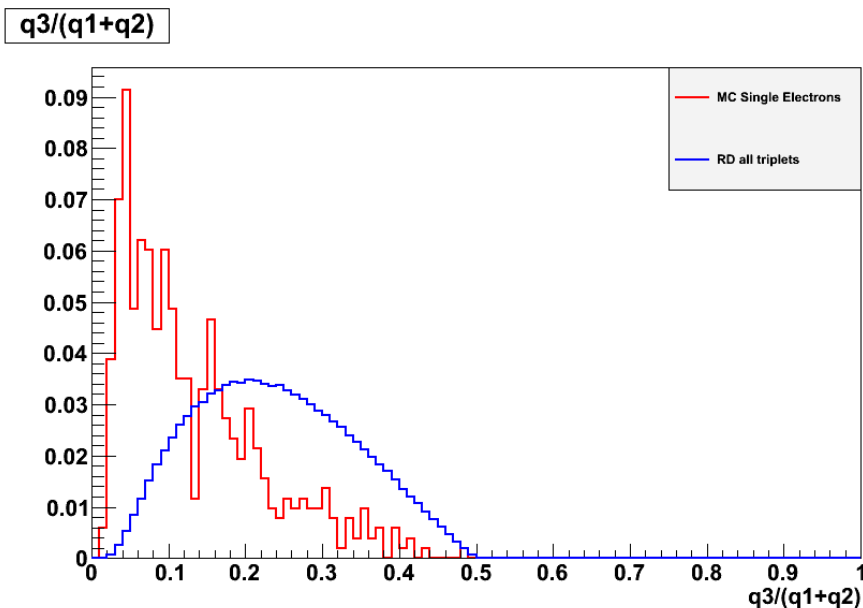
clusters {dphi!=-9999&&dz!=-9999}

htemp	
Entries	2285
Mean	7.405
RMS	1.458



Cluster shape

- Distribution of charge among triplet member pads can be used to select preclusters
 - Tried two variables $q1/(q1+q2+q3)$ and $q3/(q1+q2)$ where $q1$ to $q3$ are the charges measured in the three pads of the triplet in decreasing order
 - There seems to be some possibility to use these or similar variables but it requires serious validation of the MC response of the HBD



Summary

- A new clusterization algorithm
 - Preclusterization: all triplets, s, b, shape
 - Selection:
 - Tighter selection criteria at this step => Loss of efficiency but also more stable results in terms of cluster size and charge
 - Optimization is simple and possible
 - Merging is straight forward if selection is done well
- The geometry use inside the clusterizer is validated using single electron simulation
- How a selection criteria can be optimized is demonstrated using s/b cut and faux scintillation background
- Other potential selection parameters (shape) distribs shown.
- Things left to do:
 - Make the scintillation background more realistic
 - See the doubles responses
 - Less urgent but still important: Optimize the code itself, make it leak free